

Remarks

Claims 1-18 are pending. Applicants propose amending claims 1, 4, 5 and 9 to address a prior art rejection by eliminating a group from the definition of Ar₁ and a number of 112 issues. Claim 1 has been amended to eliminate the C₆-C₂₀aryloyl groups and to remove the reference to unsubstituted in the definition for Ar₁. Claims 4 and 5 have been amended to rearrange a substituent condition and to remove the reference to unsubstituted for the definition of Ar₁. Claim 9 has been amended to change its dependency. No new matter has been added. Applicants submit that the amendments further prosecution by addressing a number of rejections and placing the case in condition for allowance. For these reasons, Applicants submit that good cause exists to enter the amendments even though presented after final rejection.

The Examiner objects to claim 9 for failing to further limit the subject matter of a preceding claim. The Examiner believes that claim 9 should depend from claim 8, rather than claim 7. Claim 9 has been amended accordingly.

The Examiner rejects claims 1-18 under 35 U.S.C. 112(2) as being indefinite. The Examiner maintains that the definition of Ar₁ is unclear. With respect to lines 10 and 11, the Examiner proposes an additional "or", which has been adopted. Further, the definition for Ar₁ has been amended as follows: "...or C₆-C₂₀aryloyl group or with one of the carbon atoms of the aryl ring of the C₆-C₂₀aryl or C₆-C₂₀aryloyl group; or, provided that R₁ is acetyl, Ar₁ is C₃-C₉heteroaryl, which is unsubstituted or substituted 1 to 7 times by..." This wording clearly indicates that any heteroaryl, unsubstituted or substituted, is only defined for the case that R₁ is acetyl. Corresponding amendments have to be made in claims 4 and 5.

The Examiner rejects claims 1, 6, 7, 17 under the judicially created doctrine of obviousness-type double patenting over claims 6, 11, 12 and 17 of copending case 09/734,625. This rejection is provisional. Applicant shall submit a terminal disclaimer in the event this becomes the only remaining rejection, subject to a final comparison of the allowable claims herein and the scope of the copending claims. ✓

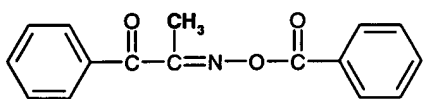
The Examiner rejects claims 1-10, 12-16 and 18 under 35 U.S.C. 102(b) as being anticipated by U.S. Pat. No. 4,282,309 ("Laridon et al."). The Examiner acknowledges that the Laridon does not exemplify the claimed compounds, but argues that the suggested substitutions can be used for anticipation. Applicants respectfully traverse this rejection.

Laridon specifically only discloses compounds corresponding to oximes with $Ar_1 = \text{aroyl}$ [$Ar_1-C(R)=N-O-R_1$]. In the present claims, Applicants have deleted "aroyl" from the definition of Ar_1 , which even further distinguishes the presently claimed compounds from the ones as specifically disclosed by Laridon. Applicants request that the Examiner reconsider and withdraw her anticipation rejection of claims 1-10, 12-16 and 18 in view of Laridon.

The Examiner rejects claim 17 under 35 U.S.C. 103 as being unpatentable over Laridon. Applicants respectfully traverse this rejection.

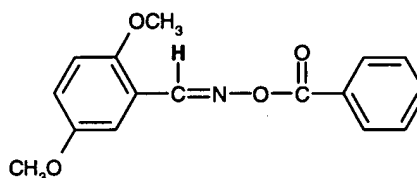
Laridon does not specifically disclose any "aldoxime" compound. In Laridon, the specific oxime ester compounds are found in col. 3, lines 20-45. In order to give a demonstration of the inventive step for the subject matter claimed in the present application, Applicants enclose a Declaration under Rule 132 of Hidetaka Oka. H. Oka compared a "ketoxime" compound as specifically disclosed in Laridon, i.e. compound A, col. 3, line 20, with two "aldoxime" compounds according to the present invention.

Compound according to
"Laridon", US Patent 4282309,
col. 3, compound (A)

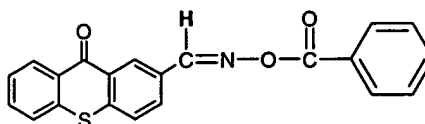


A

Compounds according to application Serial No.
09/734,635



B1



B2

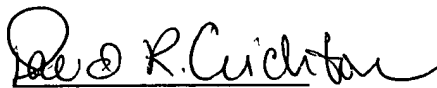
The experimental procedures are discussed more fully in the Declaration. Compounds B1 and B2 are believed to represent the closest point in view of the scope of amended claim 1. These compounds are significantly more reactive as measured using a Stouffer Wedge. The results are reproduced below show an unexpected superiority of the presently claimed compounds in resist compositions.

Compound	Number of steps reproduced after exposure time of		
	40 sec.	80 sec.	160 sec.
A	2	4	6
B1	4	7	9
B2	5	6	8

The prior art does not suggest that the selected aldoxime-type compounds would exhibit such a degree of improved performance. Thus, an inventive step is clearly given for the claimed subject matter. Applicants request that the Examiner reconsider and withdraw her obviousness rejection of claim 17 in view of Laridon.

Applicants submit that the present application is now in condition for allowance. In the event that minor amendments will further prosecution, Applicants request that the Examiner contact the undersigned representative.

Respectfully submitted,

A handwritten signature in black ink, appearing to read "David R. Crichton", written over a horizontal line.

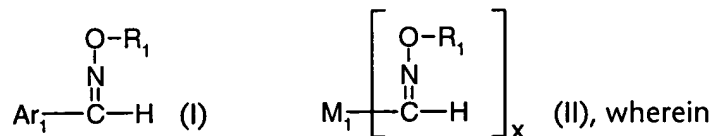
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Encl.: Declaration Under Rule 1.132, dated 18 November 2002, by H. Oka

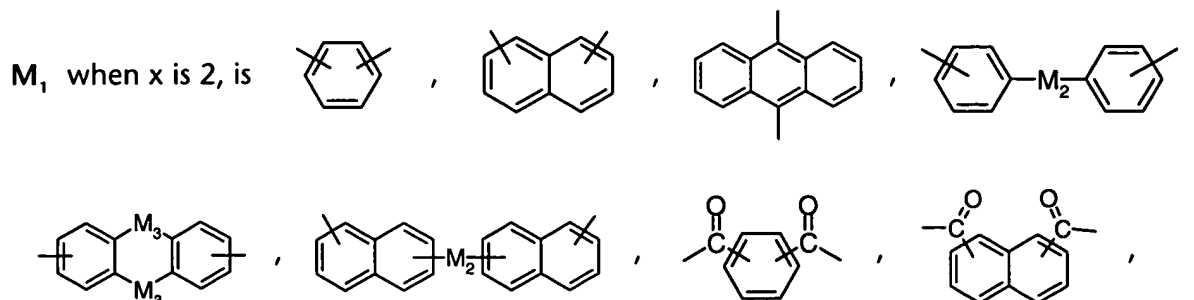
Amended Claims with underlining and bracketing

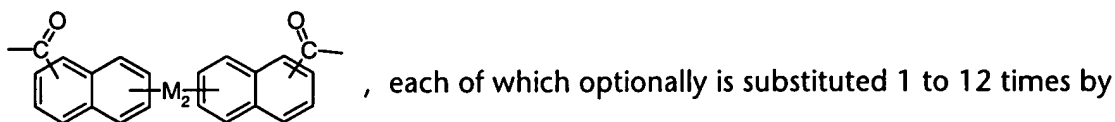
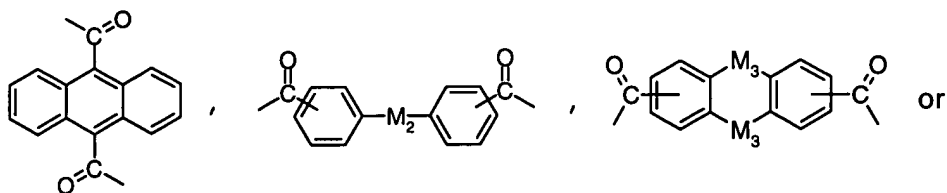
1. (amended) Alkaline developable, photosensitive composition comprising
(A) at least one alkaline soluble binder resin, prepolymer or monomer component;
(B) at least one compound of formula I or II



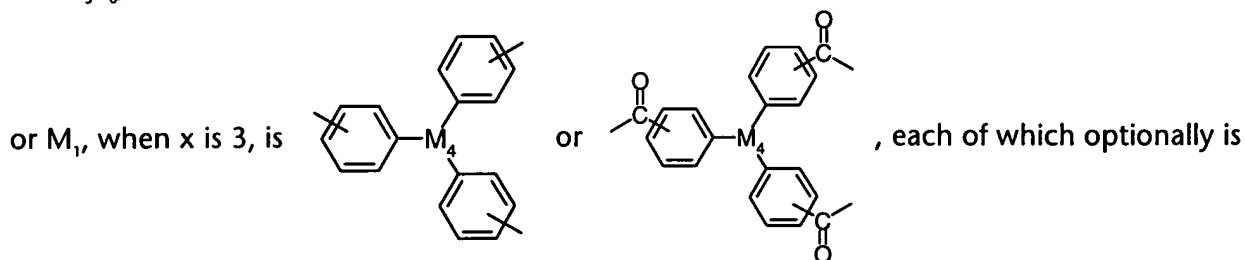
R_1 is $\text{C}_4\text{-C}_{20}$ cycloalkanoyl, $\text{C}_3\text{-C}_{12}$ alkenoyl; $\text{C}_1\text{-C}_{20}$ alkanoyl which is unsubstituted or substituted by one or more halogen, CN or phenyl; or R_1 is benzoyl which is unsubstituted or substituted by one or more $\text{C}_1\text{-C}_6$ alkyl, halogen, CN, OR_3 , SR_4 or NR_5R_6 ; or R_1 is $\text{C}_2\text{-C}_{12}$ alkoxycarbonyl or benzyloxycarbonyl; or phenoxycarbonyl which is unsubstituted or substituted by one or more $\text{C}_1\text{-C}_6$ alkyl or halogen;

Ar_1 is $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl, both radicals are unsubstituted or which is substituted 1 to 12 times by halogen, $\text{C}_1\text{-C}_{20}$ alkyl, benzyl, $\text{C}_1\text{-C}_{20}$ alkanoyl, or $\text{C}_3\text{-C}_8$ cycloalkyl; or said $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl is substituted by phenyl or benzoyl each of which optionally is substituted by one or more OR_3 , SR_4 or NR_5R_6 ; or said $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl is substituted by $\text{C}_2\text{-C}_{12}$ alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more hydroxyl groups; or said $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl is substituted by phenoxycarbonyl, OR_3 , SR_4 , SOR_4 , SO_2R_4 or NR_5R_6 , wherein the substituents OR_3 , SR_4 or NR_5R_6 optionally form 5- or 6-membered rings via the radicals R_3 , R_4 , R_5 and/or R_6 with further substituents on the aryl ring of the $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl group or with one of the carbon atoms of the aryl ring of the $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl group; or, provided that R_1 is acetyl, or Ar_1 is $\text{C}_3\text{-C}_9$ heteroaryl, provided that R_1 is acetyl, said $\text{C}_3\text{-C}_9$ heteroaryl which is unsubstituted or substituted 1 to 7 times by halogen, $\text{C}_1\text{-C}_{20}$ alkyl, benzyl, $\text{C}_1\text{-C}_{20}$ alkanoyl, or $\text{C}_3\text{-C}_8$ cycloalkyl; or said $\text{C}_3\text{-C}_9$ heteroaryl is substituted by phenyl or benzoyl, each of which optionally is substituted by one or more OR_3 , SR_4 or NR_5R_6 ; or said $\text{C}_3\text{-C}_9$ heteroaryl is substituted by $\text{C}_2\text{-C}_{12}$ alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more hydroxyl groups; or said $\text{C}_6\text{-C}_{20}$ aryl or $\text{C}_6\text{-C}_{20}$ aryloyl $\text{C}_3\text{-C}_9$ heteroaryl is substituted by phenoxycarbonyl, OR_3 , SR_4 , SOR_4 , SO_2R_4 or NR_5R_6 ;
 x is 2 or 3;





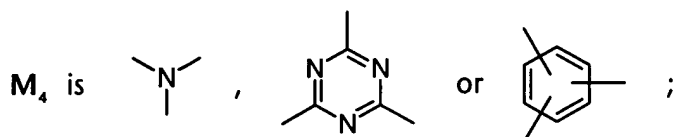
halogen, C_1 - C_{12} alkyl, C_3 - C_8 cycloalkyl, benzyl; phenyl which is unsubstituted or substituted by one or more OR_3 , SR_4 or NR_5R_6 ; benzoyl which is unsubstituted or substituted by one or more OR_3 , SR_4 or NR_5R_6 ; C_1 - C_{12} alkanoyl; C_2 - C_{12} alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more OH, phenoxy, OR_3 , SR_4 , SOR_4 , SO_2R_4 or NR_5R_6 ;



substituted 1 to 12 times by halogen, C_1 - C_{12} alkyl, C_3 - C_8 cycloalkyl; phenyl which is unsubstituted or substituted by one or more OR_3 , SR_4 or NR_5R_6 ; benzyl, benzoyl, C_1 - C_{12} alkanoyl; C_2 - C_{12} alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more hydroxyl groups, phenoxy, OR_3 , SR_4 , SOR_4 , SO_2R_4 or NR_5R_6 ;

M_2 is a direct bond, -O-, -S-, -SS-, - NR_3 -, -(CO)-, C_1 - C_{12} alkylene, cyclohexylene, phenylene, naphthylene, -(CO)O-(C_2 - C_{12} alkylene)-O(CO)-, -(CO)O-(CH_2CH_2O) $_n$ -(CO)- or -(CO)-(C₂- C_{12} alkylene)-(CO)-; or M_2 is C_4 - C_{12} alkylene or C_4 - C_{12} alkylenedioxy-, each of which is optionally interrupted by 1 to 5 -O-, -S- and/or - NR_3 -;

M_3 is a direct bond, - CH_2 -, -O-, -S-, -SS-, - NR_3 - or -(CO)-;



R_3 is hydrogen or C_1 - C_{20} alkyl; or R_3 is C_2 - C_{12} alkyl which is substituted by -OH, -SH, -CN, C_3 - C_6 alkenoxy, - OCH_2CH_2CN , - $OCH_2CH_2(CO)O(C_1$ - C_4 alkyl), -O(CO)- C_1 - C_4 alkyl, -O(CO)-phenyl, -(CO)OH, -(CO)O(C_1 - C_4 alkyl), -N(C_1 - C_4 alkyl) $_2$, -N(CH_2CH_2OH) $_2$, -N[CH_2CH_2O -(CO)- C_1 - C_4 alkyl] $_2$ or morpholinyl; or R_3 is C_2 - C_{12} alkyl which is interrupted by one or more -O-; or R_3 is -(CH_2CH_2O) $_n$ H, -(CH_2CH_2O) $_n$ (CO)- C_1 - C_8 alkyl, C_1 - C_8 alkanoyl, C_3 - C_{12} alkenyl, C_3 - C_6 alkenoyl, C_3 - C_8 cycloalkyl; or R_3 is benzoyl which is unsubstituted or substituted by one or more C_1 - C_6 alkyl, halogen, -OH or C_1 - C_4 alkoxy; or R_3 is phenyl or naphthyl each of which is unsubstituted or substituted by halogen, -OH, C_1 - C_{12} alkyl, C_1 - C_{12} alkoxy, phenyl- C_1 - C_3 -alkoxy, phenoxy, C_1 -

C₁₂alkylsulfanyl, phenylsulfanyl, -N(C₁-C₁₂alkyl)₂, diphenylamino or -(CO)R₇; or R₃ is phenyl-C₁-C₃alkyl, or Si(C₁-C₆alkyl)_r(phenyl)_{3-r};

r is 0, 1, 2 or 3;

n is 1 to 20;

R₄ is hydrogen, C₁-C₂₀alkyl, C₃-C₁₂alkenyl, C₃-C₈cycloalkyl, phenyl-C₁-C₃alkyl; C₂-C₈alkyl which is substituted by -OH, -SH, -CN, C₃-C₆alkenoxy, -OCH₂CH₂CN, -OCH₂CH₂(CO)O(C₁-C₄alkyl), -O(CO)-C₁-C₄alkyl, -O(CO)-phenyl, -(CO)OH or -(CO)O(C₁-C₄alkyl); or R₄ is C₂-C₁₂alkyl which is interrupted by one or more -O- or -S-; or R₄ is -(CH₂CH₂O)_{n+1}H, -(CH₂CH₂O)_n(CO)-C₁-C₈alkyl, C₂-C₈alkanoyl, C₃-C₁₂alkenyl, C₃-C₆alkenoyl; or R₄ is phenyl or naphthyl each of which is unsubstituted or substituted by halogen, C₁-C₁₂alkyl, C₁-C₁₂alkoxy or -(CO)R₇;

R₅ and R₆ independently of each other are hydrogen, C₁-C₂₀alkyl, C₂-C₄hydroxyalkyl, C₂-C₁₀alkoxyalkyl, C₃-C₃alkenyl, C₃-C₈cycloalkyl, phenyl-C₁-C₃alkyl, C₁-C₄alkanoyl, C₃-C₁₂alkenoyl, benzoyl; or are phenyl or naphthyl each of which is unsubstituted or substituted by C₁-C₁₂alkyl or C₁-C₁₂alkoxy; or R₅ and R₆ together are C₂-C₆alkylene optionally interrupted by -O- or -NR₃- and/or optionally substituted by hydroxyl, C₁-C₄alkoxy, C₂-C₄alkanoyloxy or benzoyloxy;

R₇ is hydrogen, C₁-C₂₀alkyl; or is C₂-C₈alkyl which is substituted by halogen, phenyl, -OH, -SH, -CN, C₃-C₆alkenoxy, -OCH₂CH₂CN, -OCH₂CH₂(CO)O(C₁-C₄alkyl), -O(CO)-C₁-C₄alkyl, -O(CO)-phenyl, -(CO)OH or -(CO)O(C₁-C₄alkyl); or R₇ is C₂-C₁₂alkyl which is interrupted by one or more -O-; or R₇ is -(CH₂CH₂O)_{n+1}H, -(CH₂CH₂O)_n(CO)-C₁-C₈alkyl, C₃-C₁₂alkenyl, C₃-C₈cycloalkyl; phenyl optionally substituted by one or more halogen, -OH, C₁-C₁₂alkyl, C₁-C₁₂alkoxy, phenoxy, C₁-C₁₂alkylsulfanyl, phenylsulfanyl, -N(C₁-C₁₂alkyl)₂, or diphenylamino; and

(D) a photopolymerizable compound.

2. (amended) Photosensitive composition according to claim 1, wherein compound (A) is an oligomeric or polymeric compound.

3. Photosensitive composition according to claim 2, wherein the photopolymerizable compound (C) is a liquid.

4. (amended) Photosensitive composition according to claim 1, wherein component (B) is a compound of formula I or II, wherein


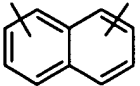
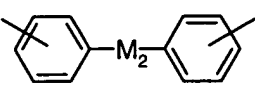
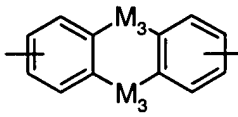
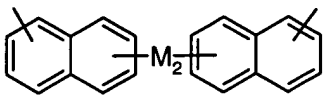
R₁ is C₂-C₆alkanoyl or C₂-C₅alkoxycarbonyl; or R₁ is benzoyl which is unsubstituted or substituted by one or more C₁-C₆alkyl or halogen;

Ar₁ is phenyl or naphthyl,

~~each of these radicals is unsubstituted or substituted 1 to 5 times by halogen, C₁-C₂₀alkyl, benzyl or C₁-C₂₀alkanoyl; or said phenyl or naphthyl is substituted by phenyl or benzoyl, each of which optionally is substituted by one or more OR₃, SR₄ or NR₅R₆; or said phenyl or naphthyl is substituted by C₂-C₁₂alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more OH; or said phenyl or naphthyl is substituted by OR₃, SR₄ or NR₅R₆, wherein the substituents OR₃, SR₄ or NR₅R₆ optionally form 5- or 6-membered rings via the radicals R₃, R₄, R₅ and/or R₆ with further substituents on the phenyl or naphthyl ring or with one of the carbon atoms of the phenyl or naphthyl ring;~~

or, provided that R₁ is acetyl, Ar₁ is furyl, pyrrolyl, thienyl, benzofuranyl, indolyl, benzothiophenyl or pyridyl, provided that R₁ is acetyl; wherein each of these radicals is unsubstituted or substituted 1 to 4 times by halogen, C₁-C₂₀alkyl, benzyl, C₃-C₈cycloalkyl,

phenyl, C₁-C₂₀alkanoyl, benzoyl, C₂-C₁₂alkoxycarbonyl, phenoxycarbonyl, OR₃, SR₄, SOR₄, SO₂R₄, or NR₅R₆;
 x is 2;

M₁ is a group , , ,  or , each of which optionally is substituted 1 to 4 times by halogen,

C₁-C₁₂alkyl, benzyl, OR₃, SR₄ or NR₅R₆; or by phenyl which is unsubstituted or substituted by one or more OR₃, SR₄ or NR₅R₆; or by benzoyl which is unsubstituted or substituted by one or more OR₃, SR₄ or NR₅R₆; or by C₁-C₁₂alkanoyl; or by C₂-C₁₂alkoxycarbonyl optionally interrupted by one or more -O- and/or optionally substituted by one or more hydroxyl groups;

M₂ is a direct bond, -O-, -S-, -SS-, -NR₃-, -(CO)-, C₁-C₁₂alkylene, phenylene, -(CO)O-(C₂-C₁₂alkylene)-O(CO)-, -(CO)O-(CH₂CH₂O)_n-(CO)- or -(CO)-(C₂-C₁₂alkylene)-(CO)-; or M₂ is C₄-C₁₂alkylene or C₄-C₁₂alkylenedioxy-, each of which is optionally interrupted by 1 to 5 -O-, -S- and/or -NR₃-;

M₃ is a direct bond, -CH₂-, -O-, -S-, -NR₃- or -(CO)-;

R₃ is hydrogen or C₁-C₂₀alkyl; or R₃ is C₂-C₁₂alkyl which is substituted by -OH, -SH, -O(CO)-C₁-C₄alkyl, -O(CO)-phenyl, -(CO)O(C₁-C₄alkyl), -N(C₁-C₄alkyl)₂, -N(CH₂CH₂OH)₂, -N[CH₂CH₂O-(CO)-C₁-C₄alkyl]₂ or morpholinyl; or R₃ is C₂-C₁₂alkyl which is interrupted by one or more -O-; or R₃ is -(CH₂CH₂O)_{n+1}H, -(CH₂CH₂O)_n(CO)-C₁-C₈alkyl, phenyl-C₁-C₃alkyl, C₂-C₈alkanoyl, C₃-C₁₂alkenyl or C₃-C₆alkenoyl; or R₃ is benzoyl which is unsubstituted or substituted by one or more C₁-C₆alkyl, halogen or C₁-C₄alkoxy; or R₃ is phenyl or naphthyl each of which is unsubstituted or substituted by halogen, C₁-C₁₂alkyl, C₁-C₁₂alkoxy, phenyl-C₁-C₃alkoxy, phenoxy, C₁-C₁₂alkylsulfanyl, phenylsulfanyl, -N(C₁-C₁₂alkyl)₂, diphenylamino or -(CO)R₇;

n is 1 to 20;

R₄ is hydrogen, C₁-C₂₀alkyl, C₃-C₁₂alkenyl, phenyl-C₁-C₃alkyl; C₂-C₈alkyl which is substituted by -OH, -SH, -O(CO)-C₁-C₄alkyl, -O(CO)-phenyl or -(CO)O(C₁-C₄alkyl); or R₄ is C₂-C₁₂alkyl which is interrupted by one or more -O- or -S-; or R₄ is -(CH₂CH₂O)_{n+1}H, -(CH₂CH₂O)_n(CO)-C₁-C₈alkyl, C₂-C₈alkanoyl, C₃-C₁₂alkenyl, C₃-C₆alkenoyl; or R₄ is phenyl or naphthyl each of which is unsubstituted or substituted by halogen, C₁-C₁₂alkyl, C₁-C₁₂alkoxy or -(CO)R₇;

R₅ and R₆ independently of each other are hydrogen, C₁-C₂₀alkyl, C₂-C₄hydroxyalkyl, C₂-C₁₀alkoxyalkyl, phenyl-C₁-C₃alkyl, C₁-C₄alkanoyl, C₃-C₁₂alkenoyl, benzoyl; or are phenyl or naphthyl each of which is unsubstituted or substituted by C₁-C₁₂alkyl or C₁-C₁₂alkoxy; or R₅ and R₆ together are C₂-C₆alkylene optionally interrupted by -O- or -NR₃- and/or optionally substituted by hydroxyl, C₁-C₄alkoxy, C₂-C₄alkanoyloxy or benzoyloxy; and

R₇ is hydrogen, C₁-C₂₀alkyl; or is C₂-C₈alkyl which is substituted by halogen, phenyl, -OH, -SH, C₃-C₆alkenoxyl, -O(CO)-C₁-C₄alkyl, -O(CO)-phenyl or -(CO)O(C₁-C₄alkyl); or R₇ is C₂-C₁₂alkyl which is interrupted by one or more -O-; or R₇ is -(CH₂CH₂O)_{n+1}H, -(CH₂CH₂O)_n(CO)-C₁-C₈alkyl or C₃-C₁₂alkenyl; or is phenyl optionally substituted by one or more halogen, C₁-C₁₂alkyl, C₁-C₁₂alkoxy, phenoxy, C₁-C₁₂alkylsulfanyl, phenylsulfanyl, -N(C₁-C₁₂alkyl)₂, or diphenylamino.

5. (amended) Photosensitive composition according to claim 1, wherein component (B) is a compound of formula I or II, wherein

R_1 is C_2 - C_4 alkanoyl;

Ar_1 is phenyl or naphthyl, each of which is ~~unsubstituted or~~ substituted by halogen, C_1 - C_8 alkyl, NR_5R_6 or OR_3 , wherein the substituents OR_3 , optionally form 5- or 6-membered rings via the radicals R_3 with further substituents on the phenyl or naphthyl ring; or, provided that R_1 is acetyl, Ar_1 is 2-furyl, 2-pyrrolyl, 2-thienyl, 3-indolyl, ~~provided that R_1 is acetyl~~;

M_1 is  ;

x is 2;

R_3 is C_1 - C_{20} alkyl; or R_3 is C_2 - C_{12} alkyl which is substituted by OH, $-O(CO)-C_1$ - C_4 alkyl, $-N(C_1$ - C_4 alkyl) $_2$, $-N(CH_2CH_2OH)_2$, $-N[CH_2CH_2O-(CO)-C_1$ - C_4 alkyl or morpholinyl; or R_3 is C_2 - C_{12} alkyl which is interrupted by one or more $-O-$; or R_3 is $-(CH_2CH_2O)_{n+1}H$ or $-(CH_2CH_2O)_n(CO)-C_1$ - C_4 alkyl;

n is 1 to 3; and

R_5 and R_6 are C_1 - C_4 alkyl.

9. Photosensitive composition according to claim 8, comprising 100 parts by weight of component (A), 0.015 to 120 parts by weight of component (B), 5 to 500 parts by weight of component (C) and 0.015 to 120 parts by weight of component (D).